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CA SUBSCRIBER PRICE	ENTRY -33.60	SESSION -33.60
=> file reg COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY 229.38	SESSION 767.82
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
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Property values tagged with IC are from the ${\tt ZIC/VINITI}$ data file provided by InfoChem.

STRUCTURE FILE UPDATES: 6 FEB 2008 HIGHEST RN 1001892-66-5 DICTIONARY FILE UPDATES: 6 FEB 2008 HIGHEST RN 1001892-66-5

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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

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=>
Uploading C:\Program Files\Stnexp\Queries\10539256IIa.str

chain nodes :
6 7 8 15 16 17 30 31
ring nodes :
1 2 3 4 5 9 10 11 12 13 14 18 19 20 21 22 23 24 25 26 27 28
29
chain bonds :
1-6 3-30 6-7 7-8 7-9 8-15 15-17 15-16 16-18 16-19 30-31
ring bonds :
1-2 1-5 2-3 3-4 4-5 9-10 9-14 10-11 11-12 12-13 13-14 18-25 18-29
19-20 19-24 20-21 21-22 22-23 23-24 25-26 26-27 27-28 28-29
exact/norm bonds :
1-2 1-5 1-6 2-3 3-4 3-30 4-5 7-8 8-15 15-17 30-31
exact bonds :
6-7 7-9 15-16 16-18 16-19
normalized bonds :

 $9-10 \quad 9-14 \quad 10-11 \quad 11-12 \quad 12-13 \quad 13-14 \quad 18-25 \quad 18-29 \quad 19-20 \quad 19-24 \quad 20-21 \quad 21-22$

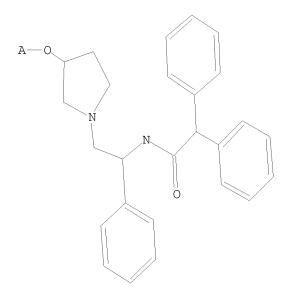
Match level:

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:CLASS 16:CLASS 17:CLASS 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom 30:CLASS 31:CLASS

L12 STRUCTURE UPLOADED

22-23 23-24 25-26 26-27 27-28 28-29

=> d 112 L12 HAS NO ANSWERS L12 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 112

SAMPLE SEARCH INITIATED 09:56:38 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 8 TO ITERATE

100.0% PROCESSED 8 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 8 TO 329
PROJECTED ANSWERS: 0 TO 0

L13 0 SEA SSS SAM L12

=> s 112 full

FULL SEARCH INITIATED 09:56:42 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 157 TO ITERATE

100.0% PROCESSED 157 ITERATIONS 2 ANSWERS

SEARCH TIME: 00.00.01

L14 2 SEA SSS FUL L12

=> file caplus

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 178.36 946.18 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE -33.60 0.00

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FILE COVERS 1907 - 7 Feb 2008 VOL 148 ISS 6 FILE LAST UPDATED: 6 Feb 2008 (20080206/ED)

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=> d ed abs ibib hitstr tot

12/02/2008,10539256IIa.trn

L15 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN

Entered STN: 01 Jul 2004

AB N-methyl-N-[(1S)-1-phenyl-2-[(3S)-3-hydroxypyrrolidin-1-yl]ethyl]-2,2-diphenylacetamide with 21 covalently bound acids and the salts,
solvates, and prodrugs thereof, were prepared for treatment of pain,
inflammation, obesity, anorexia, dysorexia, quastroparesis, dysponderosis,
etc. (no data). Thus, N-methyl-N-[(1S)-1-phenyl-2-[(3S)-3hydroxypyrrolidin-1-yl]ethyl]-2,2-diphenylacetamide hydrochloride was
warmed with Ac2O and EtSN for 2 h on a steam bath to give
N-[2-[(3S)-3-acetoxy-2-pyrrolidinyl]-(1S)-1-phenylethyl]-2,2-diphenyl-Nmethylacetamide.

ACCESSION NUMBER: 2004:525892 CAPLUS
DCCUMPNIN NUMBER: 141.89302

ACCESSION NUMBER: DOCUMENT NUMBER: TITLE: covalently 141:89002 Preparation of asimadoline derivatives with

bound acids as opiater receptor agonists
Seyfried, Christoph; Wiesner, Matthias
Merck Patent GmbH, Germany
Ger. Offen., 20 pp.
CODEN: GMXXEX
Patent
German
1 INVENTOR(S):

PATENT ASSIGNEE(S): SOURCE:

DOCUMENT TYPE: LANGUAGE: FAMILY ACC. NUM. COUNT: PATENT INFORMATION:

KIND DATE AU 2003288159 A1 20040709 AU 2003-288159 20031125
EP 1572640 A1 20050914 EP 2003-780043 20031125
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NI, SE, MC, PT,
IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, RG, CZ, EE, HU, SK
JP 200651521 T1 20060406 UP 2004-599517 20031125
US 2006122255 A1 20060608 US 2005-539256 20050616
US 2007179098 A1 20070802 US 2007-732309 20070402
RITY APPLN. INFO: DE 2002-10259245 A 20021217 PRIORITY APPLN INFO : WO 2003-EP13206 W 20031125

US 2005-539256

A1 20050616

OTHER SOURCE(S): MARPAT 141:89002

L15 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN (Continued)
IT 714237-72-6P 714237-73-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES

(claimed compound; preparation of asimadoline derivs. with covalently

acids as opiate receptor agonists) 714237-72-6 CAPLUS Benzeneacetamide, N-methyl- α -phenyl-N-[(1S)-1-phenyl-2-[(3R)-3-(sulfooxy)-1-pyrrolidinyl]ethyl]- (CA INDEX NAME)

Absolute stereochemistry.

Absolute stereochemistry.

714237-73-7 CAPLUS Benzeneacetamide, N-[(1S)-2-[(3R)-3-(acetyloxy)-1-pyrrolidiny1]-1-phenylethyl)-N-methyl-a-phenyl- (CA INDEX NAME)